

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072222\
 Data File : BG054365.D
 Acq On : 22 Jul 2022 18:38
 Operator : CG/JU
 Sample : N3835-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MH-191E

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG054365.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.568	364	371	385	rVB	293425	508730	25.92%	5.017%
2	7.172	637	644	655	rBV	334203	581703	29.64%	5.737%
3	7.989	777	783	791	rBV	57820	99019	5.05%	0.977%
4	9.152	973	981	997	rBB	201247	375185	19.12%	3.700%
5	10.797	1254	1261	1268	rBV	79870	143994	7.34%	1.420%
6	13.247	1671	1678	1690	rBB	712416	1110981	56.62%	10.957%
7	14.628	1906	1913	1919	rBV	151097	242397	12.35%	2.391%
8	16.115	2159	2166	2179	rBV2	654081	1039720	52.98%	10.255%
9	17.372	2373	2380	2384	rBV	203483	327006	16.66%	3.225%
10	17.413	2384	2387	2394	rVV	70869	101213	5.16%	0.998%
11	17.495	2396	2401	2407	rVB3	14405	27849	1.42%	0.275%
12	18.247	2523	2529	2533	rBV	16061	29441	1.50%	0.290%
13	18.294	2533	2537	2545	rVB3	21496	35376	1.80%	0.349%
14	18.441	2556	2562	2565	rBV4	19860	38475	1.96%	0.379%
15	18.741	2607	2613	2617	rBV	13743	21891	1.12%	0.216%
16	19.158	2680	2684	2690	rBV2	16550	31663	1.61%	0.312%
17	19.305	2704	2709	2715	rVB3	11613	21132	1.08%	0.208%
18	19.423	2723	2729	2735	rVB	226369	316903	16.15%	3.126%
19	19.752	2780	2785	2786	rBV	42699	50535	2.58%	0.498%
20	19.781	2786	2790	2799	rVV	199018	315175	16.06%	3.108%
21	19.851	2799	2802	2809	rVB4	14344	24409	1.24%	0.241%
22	19.981	2816	2824	2830	rBV	1410712	1962339	100.00%	19.354%
23	20.104	2839	2845	2847	rBV3	19517	34295	1.75%	0.338%
24	20.274	2864	2874	2879	rVB3	31253	76039	3.87%	0.750%
25	20.433	2896	2901	2905	rVB3	22967	38475	1.96%	0.379%
26	20.574	2921	2925	2930	rBV	17583	32299	1.65%	0.319%
27	21.032	2999	3003	3009	rBV	30755	50705	2.58%	0.500%
28	21.215	3030	3034	3037	rBV2	15201	22033	1.12%	0.217%
29	21.273	3038	3044	3047	rBV2	24101	50398	2.57%	0.497%
30	21.367	3057	3060	3068	rVB	23869	38152	1.94%	0.376%
31	21.632	3097	3105	3110	rBV3	274700	657992	33.53%	6.490%
32	21.679	3110	3113	3122	rVB	112606	186297	9.49%	1.837%
33	21.826	3135	3138	3144	rVB3	19079	29005	1.48%	0.286%
34	23.083	3346	3352	3361	rVB3	67797	146698	7.48%	1.447%
35	23.823	3469	3478	3485	rBV2	96997	339410	17.30%	3.348%
36	24.534	3592	3599	3610	rVB2	59208	170016	8.66%	1.677%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	24.681	3616	3624	3635	rVB	82179	243374	12.40%	2.400%
38	24.834	3642	3650	3659	rBV2	137900	405661	20.67%	4.001%
39	28.494	4264	4273	4289	rVB4	47181	213154	10.86%	2.102%

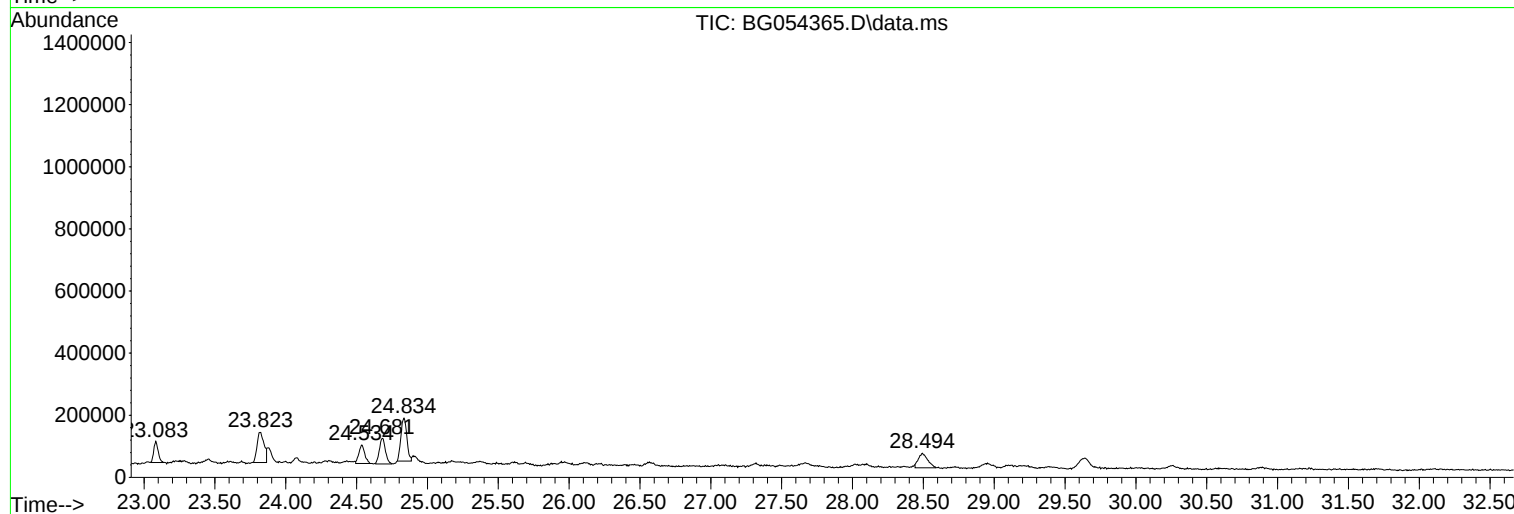
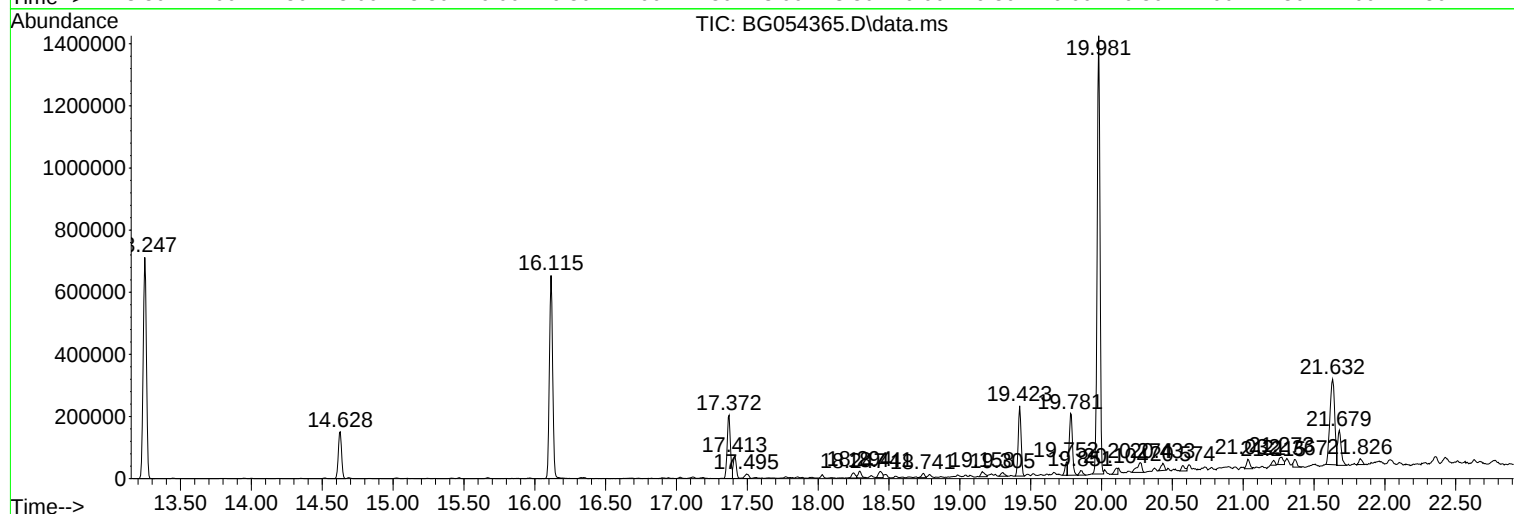
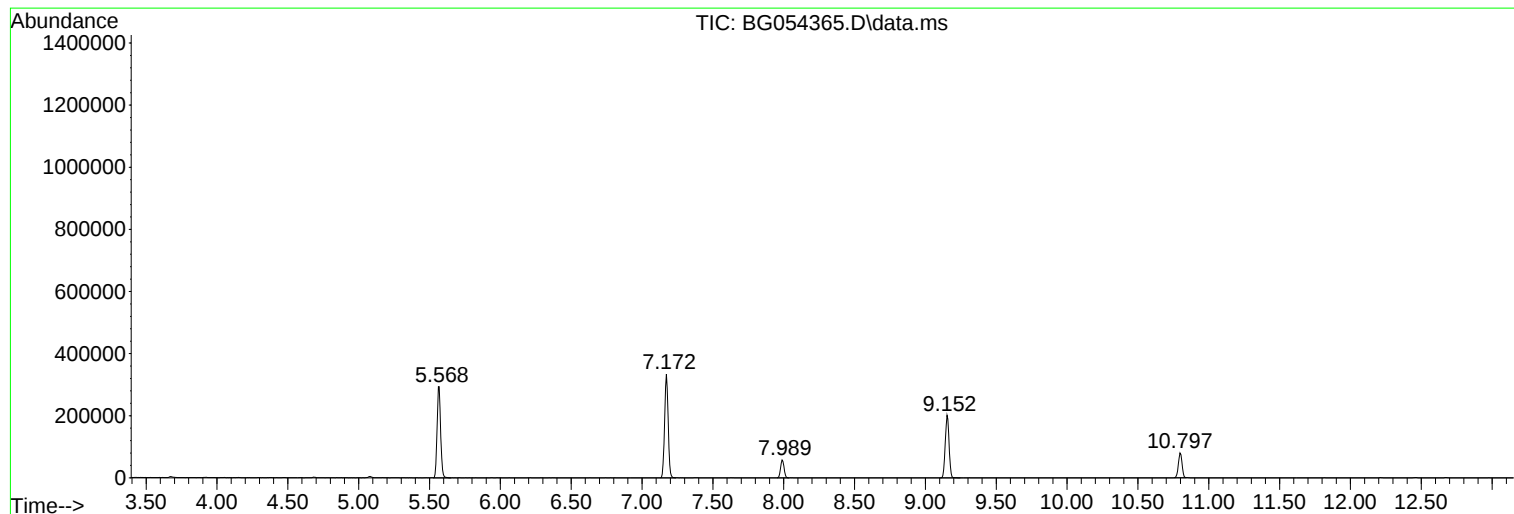
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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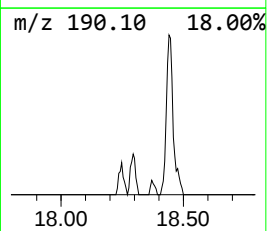
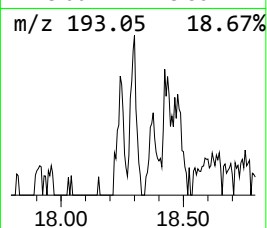
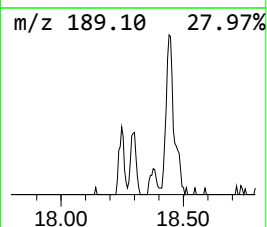
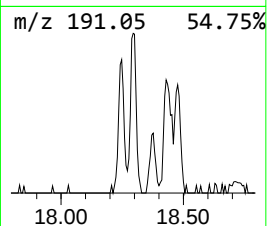
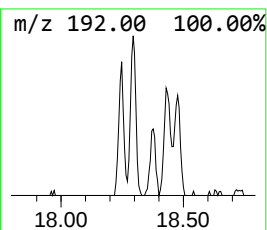
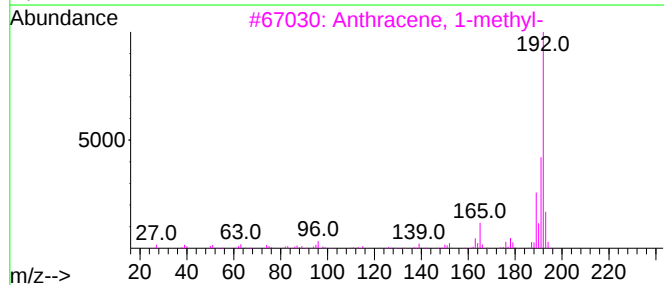
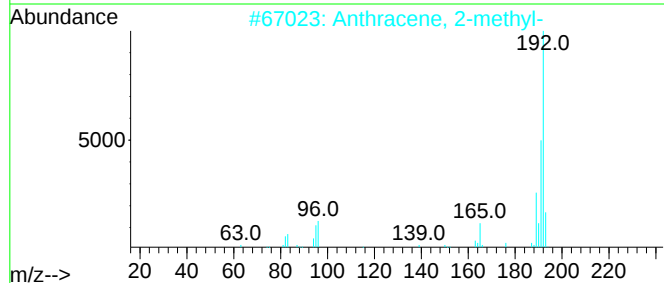
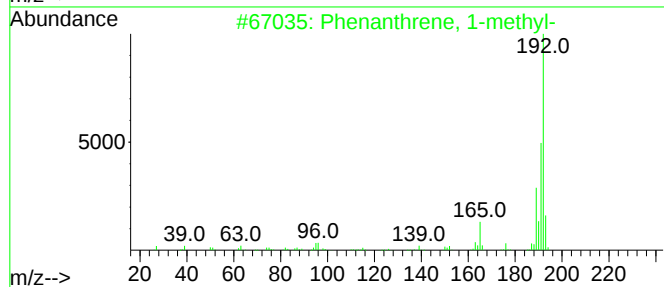
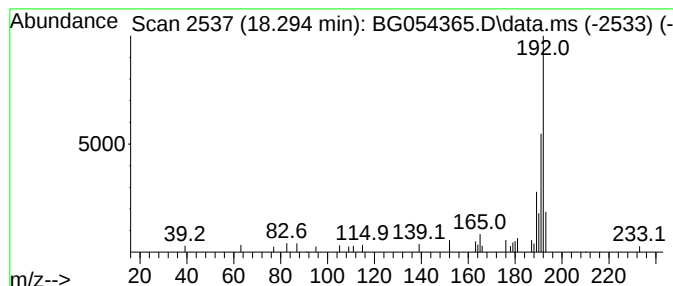
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Phenanthrene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.294	2.16 ng	35376	Phenanthrene-d10	17.372

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	81
2		Anthracene, 2-methyl-	192	C15H12	000613-12-7	81
3		Anthracene, 1-methyl-	192	C15H12	000610-48-0	81
4		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	81
5		Anthracene, 9-methyl-	192	C15H12	000779-02-2	81



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ClientSampleId :
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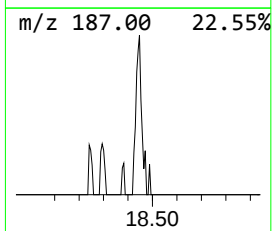
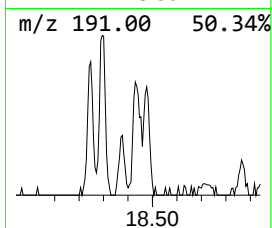
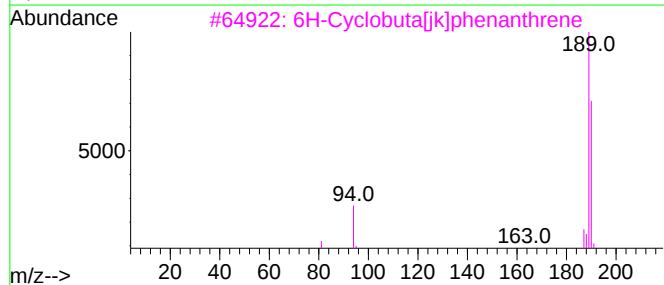
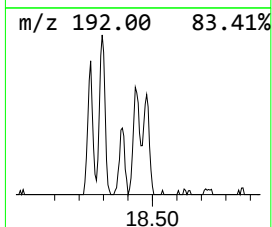
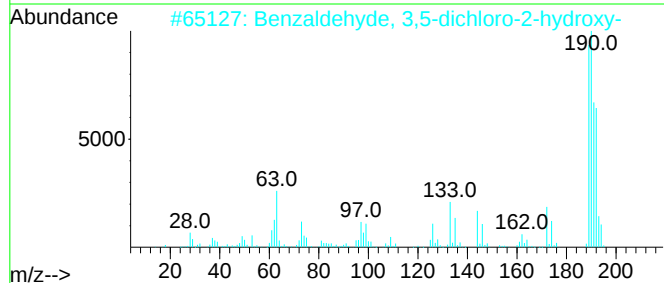
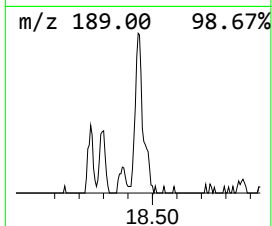
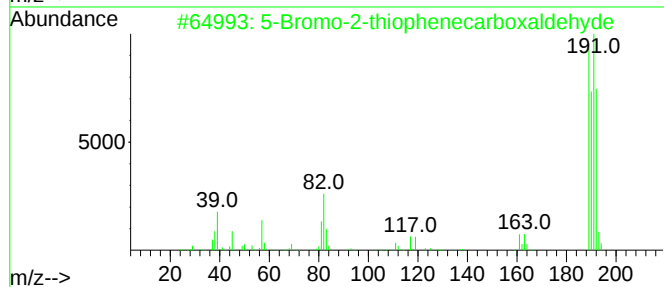
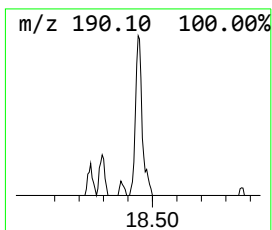
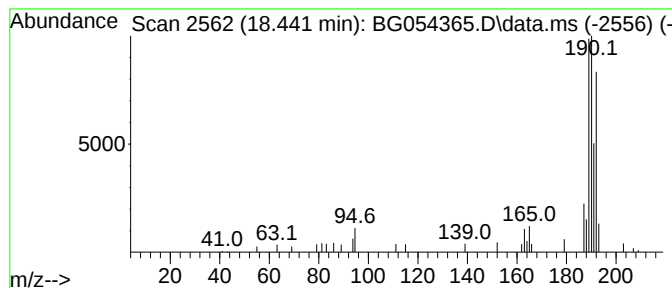
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Peak Number 2 5-Bromo-2-thiophenecarboxal... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.441	2.35 ng	38475	Phenanthrene-d10	17.372

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		5-Bromo-2-thiophenecarboxaldehyde	190	C5H3BrOS	004701-17-1	64
2		Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	58
3		6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	49
4		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	46
5		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	38



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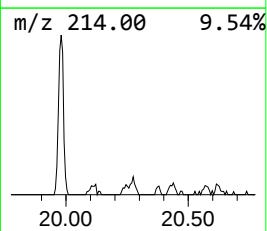
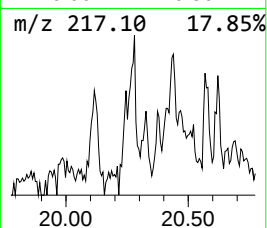
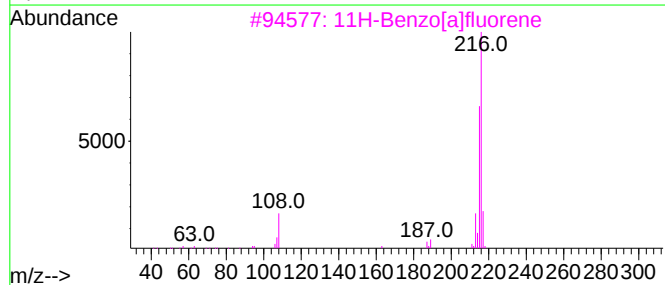
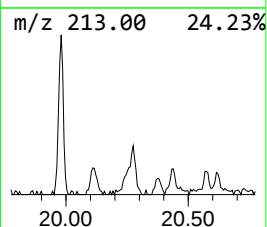
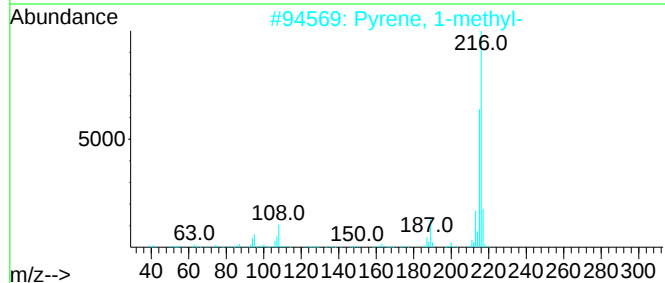
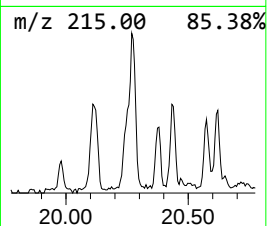
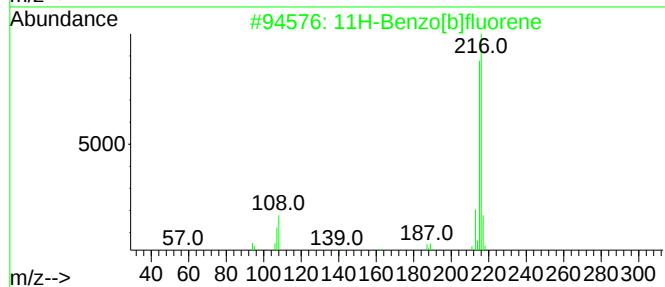
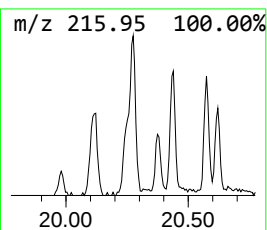
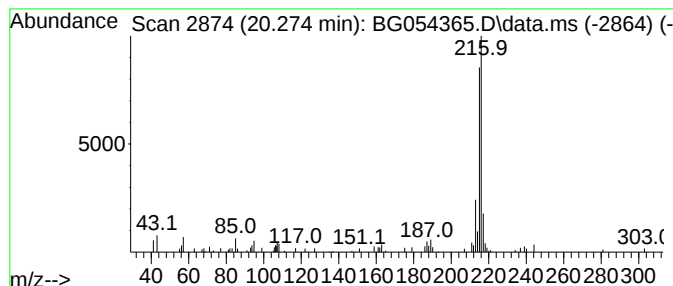
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 11H-Benzo[b]fluorene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.275	2.31 ng	76039	Chrysene-d12	21.632

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90
2			Pyrene, 1-methyl-	216	C17H12	002381-21-7	90
3			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	90
4			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	81
5			7H-Benzo[c]fluorene	216	C17H12	000205-12-9	74



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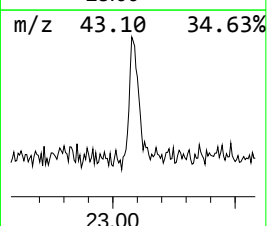
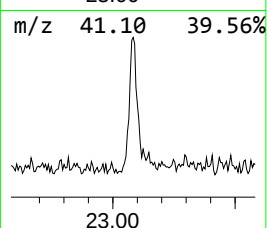
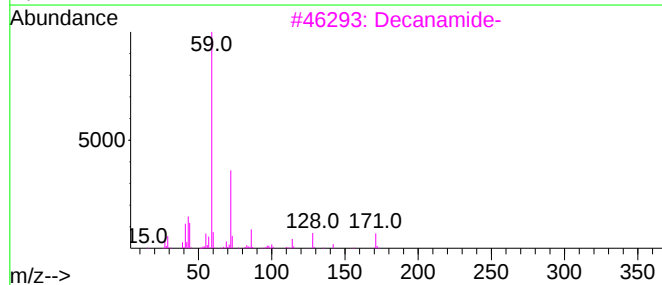
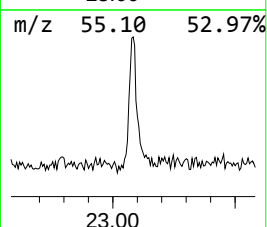
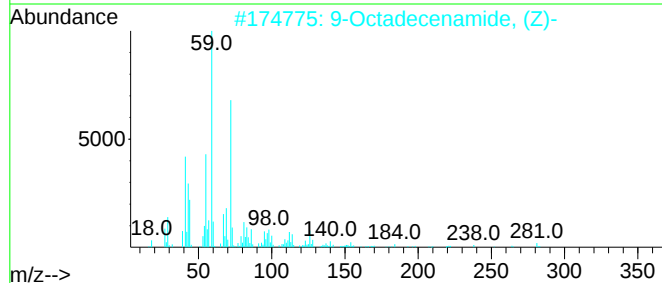
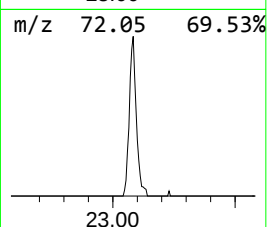
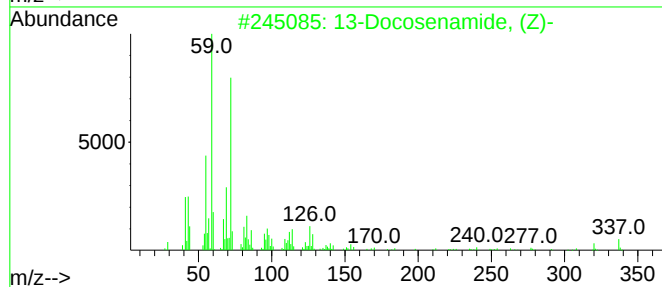
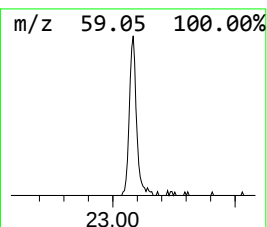
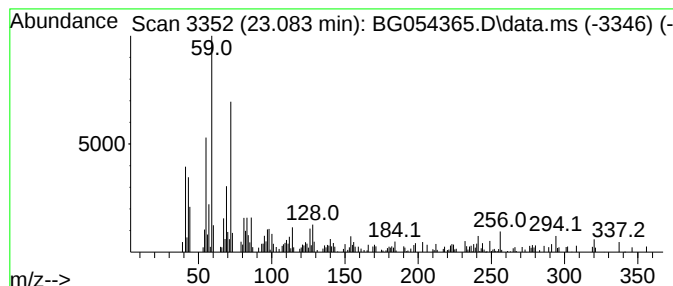
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 13-Docosenamide, (Z)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.083	4.46 ng	146698	Chrysene-d12	21.632

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	93
2		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	76
3		Decanamide-	171	C10H21NO	002319-29-1	64
4		Tetradecanamide	227	C14H29NO	000638-58-4	59
5		Palmitoleamide	253	C16H31NO	106010-22-4	59



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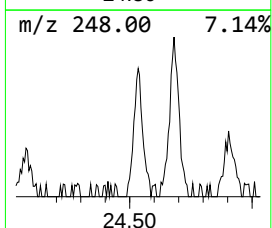
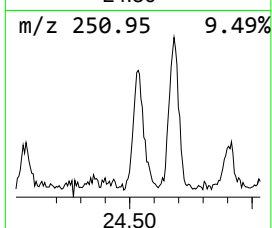
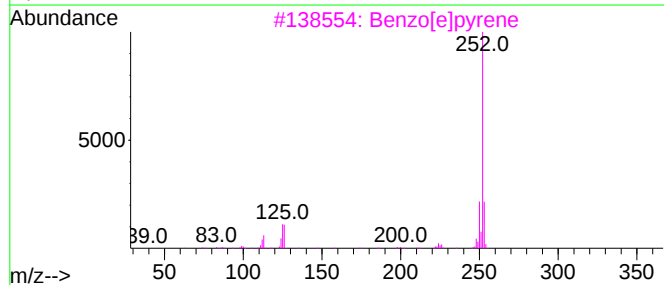
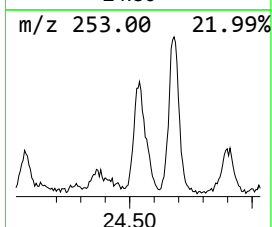
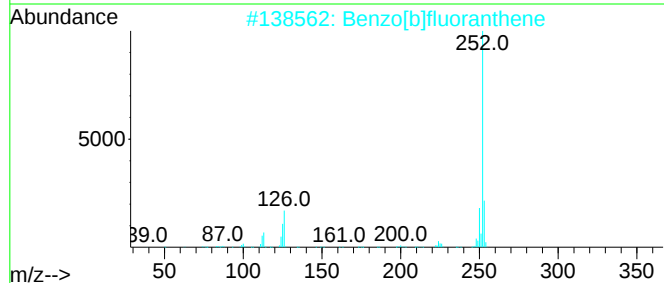
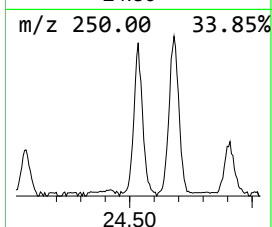
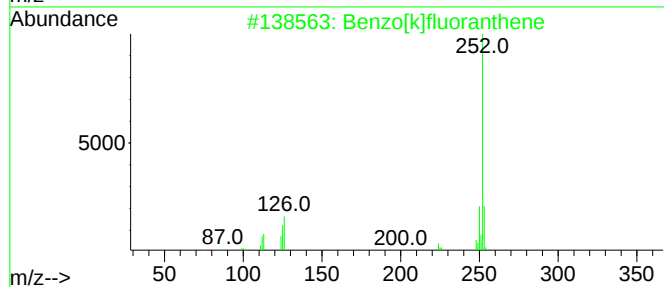
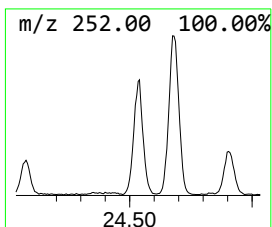
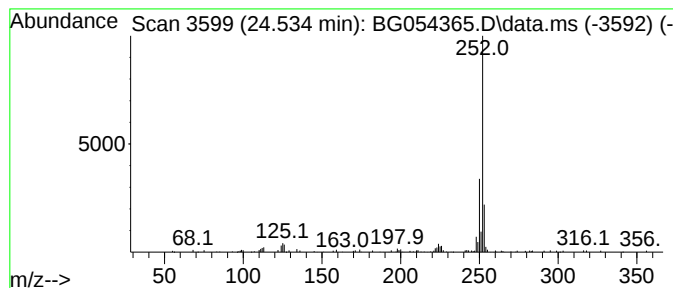
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzo[e]pyrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.534	8.38 ng	170016	Perylene-d12	24.828

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[k]fluoranthene	252	C20H12	000207-08-9	90
2			Benzo[b]fluoranthene	252	C20H12	000205-99-2	90
3			Benzo[e]pyrene	252	C20H12	000192-97-2	83
4			Perylene	252	C20H12	000198-55-0	81
5			Benzo[a]pyrene	252	C20H12	000050-32-8	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072222\
Data File : BG054365.D
Acq On : 22 Jul 2022 18:38
Operator : CG/JU
Sample : N3835-01
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
MH-191E

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Phenanthrene, 1...	18.294	2.2	ng	35376	4	17.372	327006	20.0
5-Bromo-2-thiop...	18.441	2.4	ng	38475	4	17.372	327006	20.0
11H-Benzo[b]flu...	20.275	2.3	ng	76039	5	21.632	657992	20.0
13-Docosenamide...	23.083	4.5	ng	146698	5	21.632	657992	20.0
Benzo[e]pyrene	24.534	8.4	ng	170016	6	24.828	405661	20.0